AMENDMENT AND RESPONSE TO OFFICE ACTION U.S.S.N. 10/522,595

Amendment to the Claims

This listing of claims will replace all prior versions, and listings, of claims in the application.

Listing of Claims

1-46. (Cancelled)

47. (Currently amended) A compound of formula (I), or a pharmaceutically acceptable salt or prodrug thereof

$$R_1$$
 R_2
 R_3
 R_3

wherein

X and X' taken together form $-C(R_5)=N$ -;

Y is $-C(R_5)$ - and taken together with the carbon atom bearing the phenyl group forms a double bond and R_1 is absent;

Y' is $-N(R_5)$ -;

Z forms a covalent single bond between X' and Y';

 R_2 and R_4 are independently selected from hydrogen and C_{1-3} alkyl;

 R_3 is selected from C_{1-3} alkyl and $(A)_m R_{12}$;

-C(R_5)- is selected from -C(H)- and -C(C_{1-20} alkyl)-;

-N(R₅)- is selected from -N(H)- and -N(C₂₋₂₀alkyl)-; m is 0;

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R₁₂ is selected from the group consisting of OH, SH, NH₂, halo, NO₂, C(R₁₇)₃,

 $OC(R_{17})_3$ and CN; and

R₁₇ is independently selected from hydrogen and halogen; and wherein each alkyl may be optionally substituted.

48. (Currently amended) A compound of formula (I), or a pharmaceutically acceptable salt or prodrug thereof

$$R_4$$
 R_3
 R_2
 R_3

wherein

X and X' taken together form $-C(R_5)=N$ -;

Y is $-C(R_5)$ - and taken together with the carbon atom bearing the phenyl group forms a double bond and R_1 is absent;

Y' is $-N(R_5)$ -;

Z forms a covalent single bond between X' and Y';

R₂ and R₄ are independently selected from hydrogen and C₁₋₃alkyl;

 R_3 is selected from C_{1-3} alkyl and $(A)_m R_{12}$;

 $-C(R_5)$ - is selected from -C(H)- and $-C(C_{1-20}alkyl)$ -;

-N(R₅)- is selected from -N(alkyl)- wherein alkyl is selected from the group consisting of ethyl, n-propyl, tso-propyl, cyclopropyl, n-butyl, sec-butyl, t-butyl, cyclobutyl, n-pentyl, 1-methylbutyl, 2-methylbutyl, 3-methylbutyl, cyclopentyl, n-hexyl, 1-methylpentyl, 2-methylpentyl, 3-methylpentyl, 4-methylpentyl, 1-ethylbutyl, 2-ethylbutyl, 3-ethylbutyl, 1-propylpropyl, 2-propylpropyl and cyclohexyl;

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m is 0;

R₁₂ is selected from the group consisting of OH, SH, NH₂, halo, NO₂, C(R₁₇)₃,

 $OC(R_{17})_3$ and CN;

R₁₇ is independently selected from hydrogen and halogen; and wherein each alkyl may be optionally substituted.

49. (New) The compound according to claim 47 or claim 48, or a pharmaceutically acceptable salt or prodrug thereof, wherein

Y is -CH-; and

X is -CH-.

- 50. (Previously presented) The compound according to claim 47 or claim 48, or a pharmaceutically acceptable salt or prodrug thereof, wherein R_3 is $OC(R_{17})_3$.
- 51. (Previously presented) The compound according to claim 47 or claim 48, or a pharmaceutically acceptable salt or prodrug thereof, wherein R_3 is C_{1-3} alkyl.
- 52. (Previously presented) The compound according to claim 47 or claim 48, or a pharmaceutically acceptable salt or prodrug thereof, wherein R₃ is -CH₃ or -OCH₃.
- 53. (Currently amended) The compound according to claim 47 or claim 48, or a pharmaceutically acceptable salt or prodrug thereof, wherein $N(R_5)$ is -N(3-methylbutyl).
- 54. (Previously presented) A compound wherein the compound is 4-(4-methoxyphenyl)-1-(3-methylbutyl)-1*H*-pyrazole or a pharmaceutically acceptable salt or prodrug thereof.
- 55. (Previously presented) A compound wherein the compound is 1-(3-methylbutyl)-4-(4-methylphenyl)-1*H*-pyrazole or a pharmaceutically acceptable salt or prodrug thereof.
- 56. (Previously presented) A pharmaceutical composition comprising a compound according to any one of claims 47, 48, 54 or 55, and a pharmaceutically acceptable carrier, diluent or excipient.
- 57. (Previously presented) The pharmaceutical composition according to claim 56 further comprising a glucocorticoid.